Finding heaviest H-subgraphs in real weighted graphs, with applications [*]

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Abstract

For a graph G with real weights assigned to the vertices (edges), the MAX H-SUBGRAPH problem is to find an H-subgraph of G with maximum total weight, if one exists. Our main results are new strongly polynomial algorithms for the MAX H-SUBGRAPH problem. Some of our algorithms are based, in part, on fast matrix multiplication.

For vertex-weighted graphs with n vertices we solve a more general problem – the *all pairs* MAX H-SUBGRAPH problem, where the task is to find for every pair of vertices u, v, a maximum H-subgraph containing both u and v, if one exists. We obtain an $O(n^{t(\omega,h)})$ time algorithm for the *all pairs* MAX H-SUBGRAPH problem in the case where H is a fixed graph with h vertices and $\omega < 2.376$ is the exponent of matrix multiplication. The value of $t(\omega, h)$ is determined by solving a small integer program. In particular, heaviest triangles for all pairs can be found in $O(n^{2+1/(4-\omega)}) \le o(n^{2.616})$ time. For $h = 4, 5, 8$ the running time of our algorithm essentially matches that of the (unweighted) H-subgraph detection problem. Using rectangular matrix multiplication, the value of $t(\omega, h)$ can be improved; for example, the runtime for triangles becomes $O(n^{2.575})$.

We also present improved algorithms for the MAX H-SUBGRAPH problem in the edge-weighted case. In particular, we obtain an $O(m^{2-1/k} \log n)$ time algorithm for the heaviest cycle of length $2k$ or $2k - 1$ in a graph with m edges and an $O(n^3/\log n)$ time randomized algorithm for finding the heaviest cycle of any fixed length.

Our methods also yield efficient algorithms for several related problems that are faster than any previously existing algorithms. For example, we show how to find chromatic H-subgraphs in edge-colored graphs, and how to compute the most significant bits of the distance product of two real matrices, in truly sub-cubic time.

Key words. H-subgraph, matrix multiplication, weighted graph

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1   Introduction

Finding cliques or other types of subgraphs in a larger graph are classical problems in complexity theory and algorithmic combinatorics. Finding a maximum clique is NP-Hard, and also hard to approximate [Ha98]. This problem is also conjectured to be *not* fixed parameter tractable [DF95]. The problem of finding (induced) subgraphs on k vertices in an n-vertex graph has been studied extensively (see, e.g., [AYZ95, AYZ97, CN85, EG04, KKM00, NP85, PY81, YZ04]). All known algorithms for finding an induced subgraph on k vertices have running time $n^{\Theta(k)}$. Many of these algorithms use fast matrix multiplication to obtain improved exponents.

The main contribution of this paper is a set of improved algorithms for finding (induced) k-vertex subgraphs in a real vertex-weighted or edge-weighted graph. More formally, let G be a graph with real weights assigned to the vertices (edges). The weight of a subgraph of G is the sum of the weights of its vertices (edges). The MAX H-SUBGRAPH problem is to find an H-subgraph of maximum weight, if one exists. Some of our algorithms are based, in part, on *fast* matrix multiplication. In several cases, our algorithms use fast *rectangular* matrix multiplication algorithms. However, for simplicity reasons, we express most of our time bounds in terms of ω, the exponent of fast *square* matrix multiplications. The best bound currently available on ω is $\omega < 2.376$, obtained by Coppersmith and Winograd [CW90]. This is done by reducing each rectangular matrix product to a collection of smaller square matrix products. Slightly improved bounds can be obtained by using the best available rectangular matrix multiplication algorithms of Coppersmith [Cop97] and Huang and Pan [HP98]. In all of our algorithms we assume that the graphs are *undirected*, for simplicity. All of our results are applicable to directed graphs as well. Likewise, all of our results on the MAX H-SUBGRAPH problem hold for the analogous MIN H-SUBGRAPH problem. As usual, we use the *addition-comparison* model for handling real numbers. That is, real numbers are only allowed to be compared or added. In particular, our algorithms are strongly polynomial.

Our first algorithm applies to *vertex-weighted* graphs. In order to describe its complexity we need to define a small (constant size) integer optimization problem. Let $h \geq 3$ be a positive integer. The function $t(\omega, h)$ is defined by the following optimization program.

Definition 1.1

$$b_1 = \max\{b \in N \ : \ \frac{b}{4 - \omega} \ \leq \ \lfloor \frac{h - b}{2} \rfloor\}. \tag{1}$$

$$s_1 = h - b_1 + \frac{b_1}{4 - \omega}. \tag{2}$$

$$s_2(b) = \max\{h - b + \lfloor \frac{h - b}{2} \rfloor \ , \ h - (3 - \omega)\lfloor \frac{h - b}{2} \rfloor\}. \tag{3}$$

$$s_2 = \min\{s_2(b) \ : \ \lfloor \frac{h - b}{2} \rfloor \leq b \leq h - 2\}. \tag{4}$$

$$t(\omega, h) = \min\{s_1, s_2\}. \tag{5}$$

By using fast rectangular matrix multiplication, an alternative definition for $t(\omega, h)$, resulting in slightly smaller values, can be obtained (note that if $\omega = 2$, as conjectured by many researchers, fast rectangular matrix multiplication has no advantage over fast square matrix multiplication).

Theorem 1.2 *Let H be a fixed graph with h vertices. Let $G = (V, E)$ be a graph with n vertices, and $w : V \to \Re$ a weight function. For every pair of vertices $u, v \in V$, an induced H-subgraph of G containing u and v of maximum weight (if one exists), can be found in $O(n^{t(\omega, h)})$ time. In particular, the MAX H-subgraph problem can be solved in $O(n^{t(\omega, h)})$ time.*

Notice that Theorem 1.2 solves, in fact, a more general problem, the *All-Pairs* MAX H-Subgraph problem. It is easy to establish some small values of $t(\omega, h)$ directly. For $h = 3$ we have $t(\omega, 3) = 2 + 1/(4 - \omega) < 2.616$ by taking $b_1 = 1$ in (1). Using fast rectangular matrix multiplication this can be improved to 2.575. In particular, for each pair of vertices, a triangle of maximum weight containing them (if one exists) can be found in $o(n^{2.575})$ time. This should be compared to the well-known $O(n^\omega) \le o(n^{2.376})$ time algorithm for detecting a triangle in an *unweighted* graph [IR78]. For $h = 4$ we have $t(\omega, 4) = \omega + 1 < 3.376$ by taking $b = 2$ in (4). Interestingly, the fastest algorithm for detecting a K_4, that uses square matrix multiplication, also runs in $O(n^{\omega+1})$ time [NP85]. The same phenomena also occurs for $h = 5$ where $t(\omega, 5) = \omega + 2 < 4.376$ and for $h = 8$ where $t(\omega, 8) = 2\omega + 2 < 6.752$. We also note that $t(\omega, 6) = 4 + 2/(4 - \omega)$, $t(\omega, 7) = 4 + 3/(4 - \omega)$, $t(\omega, 9) = 2\omega + 3$ and $t(\omega, 10) = 6 + 4/(4 - \omega)$. However, a closed formula for $t(\omega, h)$ cannot be given. Already for $h = 11$, and for infinitely many values thereafter, $t(\omega, h)$ is only piecewise linear in ω. For example, if $7/3 \le \omega < 2.376$ then $t(\omega, 11) = 3\omega + 2$, and if $2 \le \omega \le 7/3$ then $t(\omega, 11) = 6 + 5/(4 - \omega)$. Finally, it is easy to verify that both s_1 in (2) and s_2 in (4) converge to $3h/(6 - \omega)$ as h increases. Thus, $t(\omega, h)$ converges to $3h/(6 - \omega) < 0.828h$ as h increases.

Prior to this work, the only known algorithm for MAX H-SUBGRAPH in the vertex-weighted case (moreover, the All-Pairs version of the problem) was the naïve $O(n^h)$ algorithm. In general, reductions to fast matrix multiplication tend to fail miserably in the case of real-weighted graph problems. The most prominent example of this is the famous All-Pairs Shortest Paths (APSP) problem. Seidel [Sei95] and Galil and Margalit [GM97] developed $\tilde{O}(n^\omega)$ algorithms for undirected unweighted graphs. However, for arbitrary edge weights, the best published algorithm known is a recent $O(n^3 / \log n)$ by Chan [Ch05]. When the edge weights are integers in $[-M, M]$, the problem is solvable in $\tilde{O}(Mn^\omega)$ by Shoshan and Zwick [SZ99], and $\tilde{O}(M^{0.681} n^{2.575})$ by Zwick [Zw02], respectively. Earlier, a series of papers in the 70's and 80's starting with Yuval [Yu76] attempted to speed up APSP directly using fast matrix multiplication. Unfortunately, these works require a model that allows infinite-precision operations in constant time.

A slight modification in the algorithm of Theorem 1.2, without increasing its running time by more than a logarithmic factor, can also answer the decision problem: "for every pair of vertices u, v, is there an H-subgraph containing u and v, whose weight is in the interval $[w_1, w_2]$ where $w_1 \le w_2$ are two given reals?" Another feature of Theorem 1.2 is that it makes a relatively small number of comparisons. For example, a heaviest triangle can be found by the algorithm using only $O(m + n \log n)$ comparisons, where m is the number of edges of G.

Since Theorem 1.2 is stated for induced H-subgraphs, it obviously also applies to not-necessarily induced H-subgraphs. However, the latter problem can, in some cases, be solved faster. For example, we show that the $o(n^{2.616})$ time bound for finding a heaviest triangle also holds if one searches for a heaviest H-subgraph in the case when H is the complete bipartite graph $K_{2,k}$.

Several H-subgraph detection algorithms take advantage of the fact that G may be sparse. Improving a result of Itai and Rodeh [IR78], Alon, Yuster and Zwick obtained an algorithm for detecting a triangle, expressed in terms of m [AYZ97]. The running time of their algorithm is $O(m^{2\omega/(\omega+1)}) \leq o(m^{1.41})$. This is faster than the $O(n^\omega)$ algorithm when $m = o(n^{(\omega+1)/2})$. The best known running times in terms of m for $H = K_k$ when $k \geq 4$ are given in [EG04]. Sparseness can also be used to obtain faster algorithms for the vertex-weighted MAX H-SUBGRAPH problem. We prove:

Theorem 1.3 *If $G = (V, E)$ is a graph with m edges and no isolated vertices, and $w : V \to \Re$ is a weight function, then a triangle of G with maximum weight (if one exists) can be found in $O(m^{(18-4\omega)/(13-3\omega)}) \leq o(m^{1.45})$ time.*

The proofs of Theorems 1.2 and 1.3, and some of their consequences, appear in Section 3. In Section 2 we first introduce a general method called *dominance computation*, motivated by a problem in computational geometry and introduced by Matousek in [Ma91], and show how it can be used to obtain a truly sub-cubic algorithm for the MAX K_3-SUBGRAPH problem. Although the running time we obtain using this method is slightly inferior to that of Theorem 1.2, we show that this method has other very interesting applications. In fact, we will show how to use it in order to efficiently solve a general buyer-seller problem from computational economics. Another interesting application of the method is the ability to compute the *the most significant bits* of the distance product $A \star B$ of two real matrices, in truly sub-cubic time (see the definition of distance products in the next section). Computing the distance product quickly has long been considered as the key to a truly sub-cubic APSP algorithm, since it is known that the time complexity of APSP is no worse than that of the distance product of two arbitrary $n \times n$ matrices.

We now turn to edge-weighted graphs. An $O(m^{2-1/\lceil k/2 \rceil})$ time algorithm for detecting the existence of a cycle of length k is given in [AYZ97]. A small improvement was obtained later in [YZ04]. However, the algorithms in both papers fail when applied to edge-weighted graphs. Using the *color coding* method, together with several additional ideas, we obtain a randomized $O(m^{2-1/\lceil k/2 \rceil})$ time algorithm in the edge-weighted case, and an $O(m^{2-1/\lceil k/2 \rceil} \log n)$ deterministic algorithm.

Theorem 1.4 *Let $k \geq 3$ be a fixed integer. If $G = (V, E)$ is a graph with m edges and no isolated vertices, and $w : E \to \Re$ is a weight function, then a maximum weight cycle of length k, if one exists, can be found with high probability in $O(m^{2-1/\lceil k/2 \rceil})$ time, and deterministically in $O(m^{2-1/\lceil k/2 \rceil} \log n)$ time.*

In a recent result of Chan [Ch05] it is shown that the distance product of two $n \times n$ matrices with real entries can be computed in $O(n^3/\log n)$ time (again, reals are only allowed to be compared

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or added; more recently, Y. Han announced an $O(n^3(\log \log n / \log n)^{5/4})$ time algorithm). We show how to reduce the MAX H-SUBGRAPH problem in edge-weighted graphs to the problem of computing a distance product.

Theorem 1.5 *Let H be a fixed graph with h vertices. If $G = (V, E)$ is a graph with n vertices, and $w : E \to \Re$ is a weight function, then an induced H-subgraph of G (if one exists) of maximum weight can be found in $O(n^h / \log n)$ time.*

We can strengthen the above result considerably, in the case where H is a cycle. For (not-necessarily induced) cycles of fixed length we can combine distance products with the color coding method and obtain:

Theorem 1.6 *Let k be a fixed positive integer. If $G = (V, E)$ is a graph with n vertices, and $w : E \to \Re$ is a weight function, a maximum weight cycle with k vertices (if exist) can be found, with high probability, in $O(n^3 / \log n)$ time.*

In fact, the proof of Theorem 1.6 shows that a maximum weight cycle with $k = o(\log \log n)$ vertices can be found in (randomized) sub-cubic time. Section 4 considers edge-weighted graphs and contains the algorithms proving Theorems 1.4, 1.5 and 1.6.

Finally, we consider the related problem of finding a certain chromatic H-subgraph in an edge-colored graph. We consider the two extremal chromatic cases. An H-subgraph of an edge-colored graph is called *rainbow* if all the edges have distinct colors. It is called *monochromatic* if all the edges have the same color. Many combinatorial problems are concerned with the existence of rainbow and/or monochromatic subgraphs.

We obtain a new algorithm that finds a rainbow H-subgraph, if one exists.

Theorem 1.7 *Let H be a fixed graph with $3k + j$ vertices, $j \in \{0, 1, 2\}$. If $G = (V, E)$ is a graph with n vertices, and $c : E \to C$ is an edge-coloring, then a rainbow H-subgraph of G (if one exists) can be found in $O(n^{\omega k + j} \log n)$ time.*

The running time in Theorem 1.7 matches, up to a logarithmic factor, the running time of the induced H-subgraph detection problem in (uncolored) graphs.

We obtain a new algorithm that finds a monochromatic H-subgraph, if one exists. For fixed H, the running time of our algorithm matches the running time of the (uncolored) H-subgraph detection problem, except for the case $H = K_3$.

Theorem 1.8 *Let H be a fixed connected graph with $3k + j$ vertices, $j \in \{0, 1, 2\}$. If $G = (V, E)$ is a graph with n vertices, and $c : E \to C$ is an edge-coloring, then a monochromatic H-subgraph of G (if one exists) can be found in $O(n^{\omega k + j})$ time, unless $H = K_3$. A monochromatic triangle can be found in $O(n^{(3+\omega)/2}) \le o(n^{2.688})$ time.*

The algorithms for edge-colored graphs yielding Theorems 1.7 and 1.8 appear in Section 5. The final section contains some concluding remarks and open problems.

2   Dominance computations

Given a set of points $\{v_1, \ldots, v_n\}$ in $\mathbb{R}^d$, the *dominating pairs problem* is to find all pairs of points (v_i, v_j) such that for all $k = 1, \ldots, d$, $v_i[k] \leq v_j[k]$. The key insight to our method is a connection between the problem of finding triangles and the well-known problem of computing dominating pairs in computational geometry. This connection was inspired by recent work of Chan [Ch05], who demonstrated how a $O(c^d n^{1+\varepsilon} + n^2)$ algorithm for computing dominating pairs in d dimensions can be used to solve the arbitrary APSP problem in $O(n^3 / \log n)$ time.

In particular, we use an elegant algorithm by Matousek for computing dominating pairs in n dimensions [Ma91]. Matousek's algorithm does a bit more than determine dominances — it actually computes a matrix D such that

$$D[i, j] = |\{k \mid v_i[k] \leq v_j[k]\}|.$$

We will call D the *dominance matrix* in the following.

Theorem 2.1 (Matousek [Ma91]) *Given a set S of n points in $\mathbb{R}^n$, the dominance matrix for S can be computed in $O\left(n^{\frac{3+\omega}{2}}\right)$ time.*

We outline Matousek's approach in the following paragraphs. For each coordinate $j = 1, \ldots, n$, sort the n points by coordinate j. This takes $O(n^2 \log n)$ time. Define the *jth rank of point v_i*, denoted as $r_j(v_i)$, to be the position of v_i in the sorted list for coordinate j.

For a parameter $s \in [\log n, n]$, make n/s pairs of Boolean matrices $(A_1, B_1) \ldots, (A_{n/s}, B_{n/s})$ defined as follows:

$$A_k[i, j] = 1 \iff r_j(v_i) \in [ks, ks + s),$$

$$B_k[i, j] = 1 \iff r_j(v_i) \geq ks + s.$$

Now, multiply A_k with B_k^T, obtaining a matrix C_k. Then $C_k[i, j]$ equals the number of coordinates c such that $v_i[c] \leq v_j[c]$, $r_c(v_i) \in [ks, ks + s)$, and $r_j(v_i) \geq ks + s$.

Therefore, letting

$$C = \sum_{k=1}^{n/s} C_k,$$

we have that $C[i, j]$ is the number of coordinates c such that $\lfloor r_c(v_i)/s \rfloor < \lfloor r_c(v_j)/s \rfloor$.

Suppose we compute a matrix E such that $E[i, j]$ is the number of c such that $v_i[c] \leq v_j[c]$ and $\lfloor r_c(v_i)/s \rfloor = \lfloor r_c(v_j)/s \rfloor$. Then, defining $D := C + E$, we have the desired matrix

$$D[i, j] = |\{k \mid v_i[k] \leq v_j[k]\}|.$$

To compute E, we use the n sorted lists. For an integer n, define $[n] = \{1, \ldots, n\}$. Then, for each pair $(i, j) \in [n] \times [n]$, we look up v_i's position p in the sorted list for coordinate j. By reading off the adjacent points less than v_i in this sorted list (*i.e.* the points at positions $p - 1$, $p - 2$, *etc.*), and stopping when we reach a point v_k such that $\lfloor r_j(v_k)/s \rfloor < \lfloor r_j(v_i)/s \rfloor$, we obtain the list

$v_{i_1}, \ldots, v_{i_\ell}$ of $\ell \leq s$ points such that, for all $x = 1, \ldots, \ell$, $v_{i_x}[j] \leq v_i[j]$ and $\lfloor r_j(v_i)/s \rfloor = \lfloor r_j(v_{i_x})/s \rfloor$. Finally, for each $x = 1, \ldots, \ell$, we add a 1 to $E[i_x, i]$. Assuming constant time lookups and constant time probes into a matrix, this entire process takes only $O(n^2 s)$ time.

The running time of the above procedure is $O(n^2 s + \frac{n}{s} n^\omega)$. Choosing $s = n^{\frac{\omega-1}{2}}$, the time bound becomes $O\left(n^{\frac{3+\omega}{2}}\right)$.

2.1 Finding a heaviest triangle in sub-cubic time

We first present a weakly polynomial deterministic algorithm, then a randomized strongly polynomial algorithm.

Theorem 2.2 *On graphs with integer weights, a maximum vertex-weighted triangle can be found in $O(n^{(\omega+3)/2} \cdot \log W)$ time, where W is the maximum weight of a triangle. On graphs with real weights, a maximum vertex-weighted triangle can be found in $O(n^{(\omega+3)/2} \cdot B)$ time, where B is the maximum number of bits in a weight.*

Proof: The idea is to obtain a procedure that, given a parameter K, returns an edge (i, j) from a triangle of weight at least K. Then one can binary search to find the weight of the maximum triangle, and try all possible vertices k to get the triangle itself.

We first explain the binary search. Without loss of generality, we assume that all edge weights are at least 1. Let W be the maximum weight of a triangle. Start by checking if there is a triangle of weight at least $K = 1$ (if not, there are no triangles). Then try $K = 2^i$ for increasing i, until there exists a triangle of weight 2^i but no triangle of weight 2^{i+1}. This i will be found in $O(\log W)$ steps. After this, we search on the interval $[2^i, 2^{i+1})$ for the largest K such that there is a triangle of weight K. This takes $O(\log W)$ time for integer weights, and $O(B)$ time for real weights with B bits of precision.

We now show how to return an edge from a triangle of weight at least K, for some given K. Let $V = \{1, \ldots, n\}$ be the set of vertices. For every $i \in V$, we make a point $f_i = (e(1), \ldots, e(n))$, where

$$e(j) = \begin{cases} K - w(i) & \text{if there is an edge from } i \text{ to } j \\ \infty & \text{otherwise.} \end{cases}$$

(In implementation, we can of course substitute a sufficiently large value in place of ∞.) We also make a point $g_i = (e'(1), \ldots, e'(n))$, where

$$e'(j) = \begin{cases} w(i) + w(j) & \text{if there is an edge from } i \text{ to } j \\ -\infty & \text{otherwise.} \end{cases}$$

Compute the dominance matrix $D(K)$ on the sets $\{f_i\}$ and $\{g_i\}$. For all edges (i, j) in the graph, check if there exists a k such that $f_i[k] \leq g_j[k]$. This can be done by examining the appropriate entry in $D(K)$. If such a k exists, then we know there is a vertex k such that

$$K - w(i) \leq w(j) + w(k) \implies K \leq w(i) + w(k) + w(j),$$

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that is, there exists a triangle of weight at least K using edge (i, j). Observe that the above works for both directed and undirected graphs. $\square$

In the above, the binary search over all possible weights prevents our algorithm from being strongly polynomial. We would like to have an algorithm that, in a comparison-based model, has a runtime with *no* dependence on the bit lengths of weights. Here we present a randomized algorithm that achieves this.

Theorem 2.3 *On graphs with real weights, a maximum vertex-weighted triangle can be found in $O(n^{(\omega+3)/2} \cdot \log n)$ expected worst-case time.*

We would like to somehow binary search over the collection of triangles in the graph to find the maximum. As this collection is $O(n^3)$, we would then have our strongly polynomial bound. Ideally, one would like to pick the "median" triangle from a list of all triangles, sorted by weight. But as the number of triangles can be $\Omega(n^3)$, forming this list is hopeless. Instead, we shall show how dominance computations allow us to efficiently and uniformly sample a triangle at random, whose weight is from any prescribed interval (W_1, W_2). If we pick a triangle at random and measure its weight, there is a good chance that this weight is close to the median weight. In fact, a binary search that randomly samples for a pivot can be expected to terminate in $O(\log n)$ time.

Let $W_1, W_2 \in \mathbb{R} \cup \{-\infty, \infty\}$, $W_1 < W_2$, and G be a vertex-weighted graph.

Definition 2.4 $\mathcal{C}(W_1, W_2)$ *is defined to be the collection of triangles in G whose total weight falls in the range $[W_1, W_2]$.*

Lemma 2.5 *One can sample a triangle uniformly at random from $\mathcal{C}(W_1, W_2)$, in $O(n^{(\omega+3)/2})$ time.*

Proof: From the proof of Theorem 2.2, one can compute a matrix $D(K)$ in $O(n^{(\omega+3)/2})$ time, such that $D(K)[i, j] \neq 0$ iff there is a vertex k such that (i, k) and (k, j) are edges, and $w(i) + w(j) + w(k) > K$. In fact, the i, j entry of $D(K)$ is the *number* of distinct vertices k with this property.

Similarly, one can compute matrices $E(K)$ and $L(K)$ such that $E(K)[i, j]$ and $L(K)[i, j]$ contain the number of vertices k such that (i, k) and (k, j) are edges, and $w(i) + w(j) + w(k) \leq K$ (for E) or $w(i) + w(j) + w(k) < K$ (for L). (This can be done by flipping the signs on all coordinates in the sets of points $\{f_i\}$ and $\{g_i\}$ from Theorem 2.2, then computing dominances, disallowing equalities for L.)

Therefore, if we take $F = E(W_2) - L(W_1)$, then $F[i, j]$ is the number of vertices k where there is a path from i to k to j, and $w(i) + w(j) + w(k) \in [W_1, W_2]$.

Let f be the sum of all entries $F[i, j]$. For each $(i, j) \in E$, choose (i, j) with probability $F[i, j]/f$. By the above, this step uniformly samples an edge from a random triangle. Finally, we look at the set of vertices S that are neighbors to both i and j, and pick each vertex in S with probability $\frac{1}{|S|}$. This step uniformly samples a triangle with edge (i, j). The final triangle is therefore chosen uniformly at random. $\square$

Observe that there is an interesting corollary to the above.

Corollary 2.6 *In any graph, one can sample a triangle uniformly at random in $O(n^\omega)$ time.*

Proof: (Sketch) Multiplying the adjacency matrix with itself counts the number of 2-paths from each vertex to another vertex. Therefore one can count the number of triangles and sample just as in the above. □

We are now prepared to give the strongly polynomial algorithm.

Proof of Theorem 2.3: Start by choosing a triangle t uniformly at random from all triangles. By the corollary, this is done in $O(n^\omega)$ time.

Measure the weight W of t. Determine if there is a triangle with weight in the range (W, ∞), in $O(n^{(\omega+3)/2})$ time. If not, return t. If so, randomly sample a triangle from (W, ∞), let W' be its weight, and repeat the search with W'.

It is routine to estimate the runtime of this procedure, but we include it for completeness. Let $T(n, k)$ be the expected runtime for an n vertex graph, where k is the number of triangles in the current weight range under inspection. In the worst case,

$$T(n, k) \le \frac{1}{k} \sum_{i=1}^{k-1} T(n, k-i) + c \cdot n^{(\omega+3)/2}$$

for some constant $c \ge 1$. But this means

$$T(n, k-1) \le \frac{1}{k-1} \sum_{i=1}^{k-2} T(n, k-i) + c \cdot n^{(\omega+3)/2},$$

so

$$
\begin{aligned}
T(n, k) &\le \left(\frac{1}{k} + \frac{k-1}{k} \right) \cdot T(n, k-1) \\
&\quad + \left(1 - \frac{k-1}{k} \right) cn^{(\omega+3)/2} \\
&= T(n, k-1) + \frac{c}{k} n^{(\omega+3)/2},
\end{aligned}
$$

which solves to $T(n, k) = O(n^{(\omega+3)/2} \log k)$. ◻

2.2 Most significant bits of a distance product

Let A and B be two $n \times n$ matrices with entries in $\Re \cup \infty$. The *distance product* $C = A \star B$ is an $n \times n$ matrix with $C[i, j] = \min_{k=1\ldots,n} A[i, k] + B[k, j]$. Clearly, C can be computed in $O(n^3)$ time in the addition-comparison model. In fact, Kerr [Ke70] showed that the distance product requires $\Omega(n^3)$ on a straight-line program using $+$ and min. However, Fredman showed in [Fr76] that the distance product of two square matrices of order n can be performed in $O(n^3 (\log \log n / \log n)^{1/3})$ time. Following a sequence of improvements over Fredman's result, Chan gave an $O(n^3 / \log n)$ time algorithm for distance products.

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Computing the distance product quickly has long been considered as the key to a truly sub-cubic APSP algorithm, since it is known that the time complexity of APSP is no worse than that of the distance product of two arbitrary $n \times n$ matrices. Practically all APSP algorithms with runtime of the form $O(n^\alpha)$ have, at their core, some form of distance product. Therefore, any improvement on the complexity of distance product is interesting.

Here we show that *the most significant bits of $A \star B$* can be computed in sub-cubic time, again with no exponential dependence on edge weights. In previous work, Zwick [Zw02] shows how to compute *approximate* distance products. Given any $\varepsilon > 0$, his algorithm computes distances d_{ij} such that the difference of d_{ij} and the exact value of the distance product entry is at most $O(\varepsilon)$. The running time of his algorithm is $O(\frac{W}{\varepsilon} \cdot n^\omega \log W)$. Unfortunately, guaranteeing that the distances are within ε of the right values, does not necessarily give any of the bits of the distances. Our strategy is to use dominance matrix computations.

Proposition 2.7 *Let $A, B \in (\mathbb{Z} \cup \{+\infty, -\infty\})^{n \times n}$. The k most significant bits of all entries in $A \star B$ can be determined in $O(2^k \cdot n^{\frac{3+\omega}{2}} \log n)$ time, assuming a comparison-based model.*

Proof: For a matrix M, let $M[i,:]$ be the ith row, and $M[:,j]$ be the jth column. For a constant K, define the set of vectors

$$M^L(K) := \{(M[i,1] - K, \ldots, M[i,n] - K) \mid i = 1, \ldots, n\}.$$

Also, define

$$M^R(K) := \{(-M[1,i], \ldots, -M[n,i]) \mid i = 1, \ldots, n\}.$$

Now consider the set of vectors $S(K) = A^L(K) \cup B^R(K)$. Using a dominance computation on $S(K)$, one can obtain the matrix $C(K)$ defined by

$$C(K)[i,j] := \begin{cases} 0 & \text{if } \exists k \text{ s.t. } u_i[k] < v_j[k], u_i \in A^L(K), v_j \in B^R(K) \\ 1 & \text{otherwise} \end{cases}$$

Then for any i, j,

$$\min_k \{A[i,k] + B[k,j]\} \geq K \iff C(K)[i,j] = 1.$$

Let W be the smallest power of 2 larger than $\max_{ij}\{A[i,j]\} + \max_{ij}\{B[i,j]\}$. Then $C(\frac{W}{2})$ gives the most significant bit of each entry in $A \star B$. To obtain the second most significant bit, compute $C(\frac{W}{4})$ and $C(\frac{3W}{4})$. The second bit of $(A \star B)[i,j]$ is given by the expression:

$$(\neg C(W)[i,j] \wedge C(\tfrac{3W}{4})[i,j]) \vee (\neg C(\tfrac{W}{2})[i,j] \wedge C(\tfrac{W}{4})[i,j]).$$

In general, to recover the first k bits of $(A \star B)$, one computes $C(\cdot)$ for $O(2^k)$ values of K. In particular, to obtain the ℓ-th bits, compute

$$\bigvee_{s=0}^{2^{\ell-1}-1} [\neg C(W(1 - \tfrac{s}{2^{\ell-1}})) \wedge C(W(1 - \tfrac{s}{2^{\ell-1}} - \tfrac{1}{2^\ell}))].$$

To see this, notice that for a fixed s, if (for any $i, j \in [n]$)

$$W(1 - \frac{s}{2^{\ell-1}} - \frac{1}{2^\ell}) \leq \min_k A[i,k] + B[k,j] < W(1 - \frac{s}{2^{\ell-1}}),$$

then the ℓ-th bit of $\min_k A[i,k] + B[k,j]$ must be 1, and if the ℓ-th bit of $\min_k A[i,k] + B[k,j]$ is 1, then there must exist an s with the above property.

The values for $C(W(1 - \frac{s}{2^{\ell-1}}))$ for even s are needed for computing the $(\ell-1)$-st bits, hence to compute the ℓ-th bits, at most $2^{\ell-2} + 2^{\ell-1}$ dominance computations are necessary. To obtain the first k bits of the distance product, one needs only $O(2^k)$ dominance product computations. $\square$

2.3 Buyer-Seller stable matching

We show how the "dominance-comparison" ideas can be used to improve the runtime for solving a problem arising in computational economics. In this problem, we have a set of buyers and a set of sellers. Each buyer has a set of items he wants to purchase, together with a maximum price for each item which he is willing to pay for that item. In turn, each seller has a set of items she wishes to sell, together with a reserve price for each item which she requires to be met in order for the sale to be completed. Formally:

Definition 2.8 *An* (n,k)-Buyer-Seller *instance consists of*

- *a set* $C = \{1, \ldots, k\}$ *of* commodities [1]

- *an* n-tuple *of* buyers $B = \{b_1, \ldots, b_n\}$ *where* $b_i = (B_i, p_i)$, *s.t.* $B_i \subseteq C$ *are the commodities desired by buyer* i, *and* $p_i : B_i \to \mathbb{R}^+$ *is the maximum price function for buyer* i

- *an* n-tuple *of* sellers $S = \{s_1, \ldots, s_n\}$ *where* $s_i = (S_i, v_i)$, *s.t.* $S_i \subseteq C$ *are the commodities owned by seller* i, *and* $v_i : S_i \to \mathbb{R}^+$ *is the reserve price function for seller* i

A sale transaction for an item l between a seller who owns l and a buyer who wants l can take place if the price the buyer is willing to pay is at least the reserve price the seller has for the item. Let us imagine that each buyer wants to do business with only one seller, and each seller wants to target a single buyer. Then the transaction between a buyer and a seller consists of all the items for which the buyer's maximum price meets the seller's reserve price.

Definition 2.9 *Given a buyer* (B_i, p_i) *and a seller* (S_j, v_j) *the* transaction set C_{ij} *is defined as* $C_{ij} = \{l \mid l \in B_i \cap S_j, \ p_i(l) \geq v_j(l)\}$. *Denote by* $\mathbf{C}$ *the* transaction matrix *with entries* $|C_{ij}|$.
The price *of* C_{ij} *is defined as* $P_{ij} = \sum_{l \in C_{ij}} p_i(l)$, *and the* reserve *of* C_{ij} *is defined as* $R_{ij} = \sum_{l \in C_{ij}} v_j(l)$. *Denote by* P *and* R *respectively the transaction* price *and* reserve *matrices with entries* P_{ij} *and* R_{ij}.

---

[1] We will use the words "commodities" and "items" interchangeably.

Further, we assume that every buyer i has a preference relation on the sellers j which depends entirely on P_{ij}, R_{ij} and $|C_{ij}|$. Conversely, every seller has a preference relation on the buyers determined by the same three values. More formally,

- buyer i has a (computable) preference function $f_i : \mathbb{R}^+ \times \mathbb{R}^+ \times \mathbb{Z}^+ \to \mathbb{Z}$ such that i prefers seller j to seller j' iff $f_i(P_{ij}, R_{ij}, |C_{ij}|) \geq f_i(P_{ij'}, R_{ij'}, |C_{ij'}|)$.

- Similarly, seller j has a (computable) preference function $g_j : \mathbb{R}^+ \times \mathbb{R}^+ \times \mathbb{Z}^+ \to \mathbb{Z}$ such that j prefers buyer i to buyer i' iff $g_j(P_{ij}, R_{ij}, |C_{ij}|) \geq g_j(P_{i'j}, R_{i'j}, |C_{i'j}|)$.

Ideally, each buyer wants to talk to his most preferred seller, and each seller wants to sell to her most preferred buyer. Unfortunately, this is not always possible for all buyers, even when the prices and reserves are all equal, and all preference functions equal $|C_{ij}|$. This is evidenced by the following example: Buyer 1 wants to buy item 2, buyer 2 wants to buy items 1 and 2, seller 1 has item 1, seller 2 has items 1 and 2. Here buyer 1 will not be able to get any items.

In a realistic setting, we want to find a buyer-seller matching so that there is no pair (b_i, s_j) for which b_i is not paired with s_j, such that both b_i and s_j would benefit from breaking their matches and pairing among each other. This is the *stable matching* problem, for which optimal algorithms are known when the preferences are known (*e.g.*, Gale-Shapley [GS62] can be implemented to run in $O(n^2)$). However, for large k, the major bottleneck in our setting is that of computing the preference functions of the buyers and sellers.

The obvious approach to compute P_{ij}, R_{ij} and $|C_{ij}|$ is to explicitly find the sets C_{ij}. This gives an $O(kn^2)$ algorithm to compute P_{ij}, R_{ij} and $|C_{ij}|$ for all pairs (i, j).

Let for a (computable) function f, T_f be the maximum time, over all n-bit p, r and c, needed to compute $f(p, r, c)$. Let T be the maximum time T_f, over all preference functions f_i and g_j for a buyer-seller instance. Then in time $O(kn^2 + Tn^2 + n^2 \log n)$ one can obtain for every buyer (seller) a list of the sellers (buyers) sorted by the buyer's (seller's) preference function. Exploiting fast dominance computation, we can do better.

Theorem 2.10 *The matrices P, R and $\mathbf{C}$ for an (n, k)-Buyer-Seller instance can be determined in $O(n\sqrt{kM(n,k)})$ time, where $M(n, k)$ is the time required to multiply an $n \times k$ by a $k \times n$ matrix.*

Proof: Using the dominance technique, we can compute matrix $\mathbf{C}$ as follows. For each buyer i we create a k-dimensional vector $\beta_i = \{\beta_{i1}, \ldots, \beta_{ik}\}$ so that $\beta_{ij} = p_i(j)$ if $j \in C_i$, and $\beta_{ij} = -\infty$ if $j \notin C_i$. For each seller i we create a k-dimensional vector $\sigma_i = \{\sigma_{i1}, \ldots, \sigma_{ik}\}$ so that $\sigma_{ij} = v_i(j)$ if $j \in S_i$, and $\sigma_{ij} = \infty$ if $j \notin S_i$. Computing the dominance matrix for these points computes exactly the number of items l which buyer i wants to buy, seller j wants to sell, and $p_i(\ell) \geq v_j(\ell)$.

By a modification of Matousek's algorithm for computing dominances, we can also compute the matrices P and R. We demonstrate how to find R. Recall that the dominance algorithm does a matrix multiplication $A_k \cdot B_k^T$ with entries $A_k[i, j] = 1$ iff $r_j(b_i) \in [ks, ks + s)$, and $B_k[i, j] = 1$ iff $r_j(s_i) \geq ks + s$ (using the notation from Theorem 2.1). Let B_k be the same, but redefine A_k to be

$$A_k[i, j] = \begin{cases} v_i(j) & \text{if } r_j(b_i) \in [ks, ks + s) \\ 0 & \text{otherwise} \end{cases}.$$

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Similar modifications are made to the computation of the matrix E. Instead of adding 1 to the matrix entry $E[i_x, i]$ in the step for coordinate j, we add the corresponding reserve price $v_{i_x}(j)$. Determining P can be done analogously. □

Corollary 2.11 *A buyer-seller stable matching can be determined in $O(n\sqrt{kM(n,k)} + n^2\log n + n^2T)$, where T is the maximum time to compute the preference functions of the buyers/sellers, given the buyer price and seller reserve sums for a buyer-seller pair.*

For instance, if $k = n$ and $T = O(\text{polylog } n)$, the runtime of finding a buyer-seller stable matching is $O(n^{\frac{3+\omega}{2}}) = O(n^{2.688})$.

3 Heaviest H-subgraphs of real vertex-weighted graphs

In the proof of Theorem 1.2 it would be convenient to assume that $H = K_h$ is a clique on h vertices. The proof for all other induced subgraphs with h vertices is only slightly more cumbersome, but essentially the same.

Let $G = (V, E)$ be a graph with real vertex weights, and assume $V = \{1, \ldots, n\}$. For two positive integers a, b, the *adjacency system* $A(G, a, b)$ is the 0-1 matrix defined as follows. Let S_x be the set of all $\binom{n}{x}$ x-subsets of vertices. The *weight* $w(U)$ of $U \in S_x$ is the sum of the weights of its elements. We *sort* the elements of S_x according to their weights. This requires $O(n^x\log n)$ time, assuming x is a constant. Thus, $S_x = \{U_{x,1}, \ldots, U_{x,\binom{n}{x}}\}$ where $w(U_{x,i}) \leq w(U_{x,i+1})$. The matrix $A(G, a, b)$ has its rows indexed by S_a. More precisely, the j'th row is indexed by $U_{a,j}$. The columns are indexed by S_b where the j'th column is indexed by $U_{b,j}$. We put $A(G, a, b)[U, U'] = 1$ if and only if $U \cup U'$ induces a K_{a+b} in G (this implies that $U \cap U' = \emptyset$), otherwise $A(G, a, b)[U, U'] = 0$. Notice that the construction of $A(G, a, b)$ requires $O(n^{a+b})$ time.

For positive integers a, b, c, so that $a + b + c = h$, consider the Boolean product $A(G, a, b, c) = A(G, a, b) \times A(G, b, c)$. For $U \in S_a$ and $U' \in S_c$ for which $A(G, a, b, c)[U, U'] = 1$, define their *maximum witness* $\delta(U, U')$ to be the maximal element $U'' \in S_b$ for which $A(G, a, b)[U, U''] = 1$ and also $A(G, b, c)[U'', U'] = 1$. For each $U \in S_a$ and $U' \in S_c$ with $A(G, a, b, c)[U, U'] = 1$ and with $U \cup U'$ inducing a K_{a+c}, if $U'' = \delta(U, U')$ then $U \cup U' \cup U''$ induces a K_h in G whose weight is the largest of all the K_h copies of G that contain $U \cup U'$. This follows from the fact that S_b is sorted. Thus, by computing the maximum witnesses of all plausible pairs $U \in S_a$ and $U' \in S_c$ we can find, for each pair of vertices, a K_h in G with maximum weight containing them, if such a K_h exists, or else determine that no K_h-subgraph contains the pair.

Let $A = A_{n_1 \times n_2}$ and $B = B_{n_2 \times n_3}$ be two 0-1 matrices. The *maximum witness matrix* of AB is the matrix $W = W_{n_1 \times n_3}$ defined as follows.

$$W[i,j] := \begin{cases} \text{the maximum } k \text{ s.t. } A[i,k] = B[k,j] = 1 & \text{if } (AB)[i,j] \neq 0, \\ 0 & \text{otherwise.} \end{cases}$$

Let $f(n_1, n_2, n_3)$ be the time required to compute the maximum witness matrix of the product of an $n_1 \times n_2$ matrix by an $n_2 \times n_3$ matrix. Let $h \geq 3$ be a fixed positive integer. For all possible choices of positive integers a, b, c with $a + b + c = h$ denote

$$f(h, n) = \min_{a+b+c=h} f(n^a, n^b, n^c).$$

Clearly, the time to sort S_b and to construct $A(G, a, b)$ and $A(G, b, c)$ is overwhelmed by $f(n^a, n^b, n^c)$. It follows from the above discussion that:

Lemma 3.1 *Let $h \geq 3$ be a fixed positive integer and let $G = (V, E)$ be a graph with n vertices, each having a real weight. For all pairs of vertices $u, v \in V$, an induced H-subgraph of G containing u and v of maximum weight (if one exists), can be found in $O(f(h, n))$ time. Furthermore, if $f(n^a, n^b, n^c) = f(h, n)$ then the number of comparisons needed to find a maximum weight K_h is $O(n^b \log n + z(G, a + c))$ where $z(G, a + c)$ is the number of K_{a+c} in G.*

In fact, if $b \geq 2$, the number of comparisons in Lemma 3.1 can be reduced to only $O(n^b + z(G, a+c))$. Sorting S_b reduces to sorting the sums $X + X + \ldots + X$ (X repeated b times) of an n-element set of reals X. Fredman showed in [Fr76a] that this can be achieved with only $O(n^b)$ comparisons.

A simple randomized algorithm for computing (not necessarily maximum) witnesses for Boolean matrix multiplication, in essentially the same time required to perform the product, is given by Seidel [Sei95]. His algorithm was derandomized by Alon and Naor [AN96]. An alternative, somewhat slower deterministic algorithm was given by Galil and Margalit [GM93]. However, computing the matrix of maximum witnesses seems to be a more difficult problem. Improving an earlier algorithm of Bender et al. [BFPSS05], Kowaluk and Lingas [KL05] show that $f(3, n) = O(n^{2+1/(4-\omega)}) \leq o(n^{2.616})$. This already yields the case $h = 3$ in Theorem 1.2. We will need to extend and generalize the method from [KL05] in order to obtain upper bounds for $f(h, n)$. Our extension will enable us to answer more general queries such as "is there a K_h whose weight is within a given weight interval?"

Proof of Theorem 1.2: Let $h \geq 3$ be a fixed integer. Suppose a, b, c are three positive integers with $a + b + c = h$ and suppose that $0 < \mu \leq b$ is a real parameter. For two 0-1 matrices $A = A_{n^a \times n^b}$ and $B = B_{n^b \times n^c}$ the μ-*split* of A and B is obtained by splitting the columns of A and the rows of B into consecutive parts of size $\lceil n^\mu \rceil$ or $\lfloor n^\mu \rfloor$ each. In the sequel we ignore floors and ceilings whenever it does not affect the asymptotic nature of our results. This defines a partition of A into $p = n^{b-\mu}$ rectangular matrices $A_1, \ldots, A_p$, each with n^a rows and n^μ columns, and a partition of B into p rectangular matrices $B_1, \ldots, B_p$, each with n^μ rows and n^c columns. Let $C_i = A_i B_i$ for $i = 1, \ldots, p$. Notice that each element of C_i is a nonnegative integer of value at most n^μ and that $AB = \sum_{i=1}^p C_i$. Given the C_i, the maximum witness matrix W of the product AB can be computed as follows. To determine $W[i, j]$ we look for the maximum index r for which $C_r[i, j] \neq 0$. If no such r exists, then $W[i, j] = 0$; otherwise, having found r, we now look for the maximal index k so that $A_r[i, k] = A_r[k, j] = 1$. Having found k we clearly have $W[i, j] = (r - 1)n^\mu + k$.

We now determine a choice of parameters a, b, c, μ so that the time to compute $C_1, \ldots, C_p$ and the time to compute the maximum witnesses matrix W, is $O(n^{t(\omega, h)})$. By Lemma 3.1, this suffices

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in order to prove the theorem. We will only consider $\mu \leq \min\{a, b, c\}$. Taking larger values of μ results in worse running times. The rectangular product C_i can be computed by performing $O(n^{a-\mu}n^{c-\mu})$ products of square matrices of order n^μ. Thus, the time required to compute C_i is

$$O(n^{a-\mu}n^{c-\mu}n^{\omega\mu}) = O(n^{a+c+(\omega-2)\mu}).$$

Since there are p such products, and since each of the n^{a+c} witnesses can be computed in $O(p+n^\mu)$ time, the overall running time is

$$O(pn^{a+c+(\omega-2)\mu} + n^{a+c}(p + n^\mu)) = O(n^{h-(3-\omega)\mu} + n^{h-\mu} + n^{h-b+\mu})$$

$$= O(n^{h-(3-\omega)\mu} + n^{h-b+\mu}). \tag{6}$$

Optimizing on μ we get $\mu = b/(4-\omega)$. Thus, if, indeed, $b/(4-\omega) \leq \min\{a, c\}$ then the time needed to find W is $O(n^{h-b+b/(4-\omega)})$. Of course, we would like to take b as large as possible under these constraints. Let, therefore, b_1 be the largest integer b so that $b/(4-\omega) \leq \lfloor(h-b)/2\rfloor$. For such a b_1 we can take $a = \lfloor(h-b_1)/2\rfloor$ and $c = \lceil(h-b_1)/2\rceil$ and, indeed, $\mu \leq \min\{a, c\}$. Thus, (6) gives that the running time to compute W is

$$O(n^{h-b_1+b_1/(4-\omega)}).$$

This justifies s_1 appearing in (2) in the definition of $t(\omega, h)$. There may be cases where we can do better, whenever $b/(4-\omega) > \min\{a, c\}$. We shall only consider the cases where $a = \mu = \lfloor(h-b)/2\rfloor \leq b$ (other cases result in worse running times). In this case $c = \lceil(h-b)/2\rceil$ and, using (6), the running time is
$$O(n^{h-(3-\omega)\lfloor\frac{h-b}{2}\rfloor} + n^{h-b+\lfloor\frac{h-b}{2}\rfloor}).$$

This justifies s_2 appearing in (4) in the definition of $t(\omega, h)$. Since $t(\omega, h) = \min\{s_1, s_2\}$ we have proved that W can be computed in $O(n^{t(\omega, h)})$ time. $\square$

As can be seen from Lemma 3.1 and the remark following it, the number of comparisons that the algorithm performs is relatively small. For example, in the case $h = 3$ we have $a = b = c = 1$ and hence the number of comparisons is $O(n \log n + m)$. In all the three cases $h = 4, 5, 6$ the value $b = 2$ yields $t(\omega, h)$. Hence, the number of comparisons is $O(n^2)$ for $h = 4$, $O(n^2 + mn)$ for $h = 5$ and $O(n^2 + m^2)$ for $h = 6$.

Suppose $w : \{1, \ldots, n^b\} \to \Re$ so that $w(k) \leq w(k+1)$. The use of the μ-split in the proof of Theorem 1.2 enables us to determine, for each i, j and for a real interval $I(i, j)$, whether or not there exists an index k so that $A[i, k] = B[k, j] = 1$ and $w(k) \in I(i, j)$. This is done by performing a binary search within the $p = n^{b-\mu}$ matrices $C_i, \ldots, C_p$. The running time in (6) only increases by a $\log n$ factor. We therefore obtain the following corollary.

Corollary 3.2 *Let H be a fixed graph with h vertices, and let $I \subset \Re$. If $G = (V, E)$ is a graph with n vertices, and $w : V \to \Re$ is a weight function, then, deciding whether G contains an induced H-subgraph with total weight in I can be done $O(n^{t(\omega, h)} \log n)$ time.*

Proof of Theorem 1.3: We partition the vertex set V into two parts $V = X \cup Y$ according to a parameter Δ. The vertices in X have degree at most Δ. The vertices in Y have degree larger than Δ. Notice that $|Y| < 2m/\Delta$. In $O(m\Delta)$ time we can scan all triangles that contain a vertex from X. In particular, we can find a heaviest triangle containing a vertex from X. By Theorem 1.2, a heaviest triangle induced by Y can be found in $O((m/\Delta)^{t(\omega,3)}) = O((m/\Delta)^{2+1/(4-\omega)})$ time. Therefore, a heaviest triangle in G can be found in

$$O\left(m\Delta + \left(\frac{m}{\Delta}\right)^{2+1/(4-\omega)}\right)$$

time. By choosing $\Delta = m^{(5-\omega)/(13-3\omega)}$ the result follows. $\square$

The results in Theorems 1.2 and 1.3 are useful not only for real vertex weights, but also when the weights are large integers. Consider, for example, the graph parameter $\beta(G, H)$, the H *edge-covering number* of G. We define $\beta(G, H) = 0$ if G has no H-subgraph, otherwise $\beta(G, H)$ is the maximum number of edges incident with an H-subgraph of G. To determine $\beta(G, K_k)$ we assign to each vertex a weight equal to its degree. We now use the algorithm of Theorem 1.2 to find the heaviest K_k. If the weight of the heaviest K_k is w, then $\beta(G, K_k) = w - \binom{k}{2}$. In particular, $\beta(G, K_k)$ can be computed in $O(n^{t(\omega,k)})$ time.

Finally, we note that Theorems 1.2 and 1.3 apply also when the weight of an H-subgraph is not necessarily defined as the sum of the weights of its vertices. Suppose that the weight of a triangle (x, y, z) is defined by a function $f(x, y, z)$ that is monotone in each variable separately. For example, we may consider $f(x, y, z) = xyz$, $f(x, y, z) = xy + xz + yz$ etc. Assuming that $f(x, y, z)$ can be computed in constant time given x, y, z, it is easy to modify Theorems 1.2 and 1.3 to find a triangle whose weight is maximum with respect to f in $O(n^{2+1/(4-\omega)})$ time and $O(m^{(18-4\omega)/(13-3\omega)})$ time, respectively.

We conclude this section with the following proposition.

Proposition 3.3 *If $G = (V, E)$ is a graph with n vertices, and $w : V \to \Re$ is a weight function, then a (not necessarily induced) maximum weight $K_{2,k}$-subgraph can be found in $O(n^{2+1/(4-\omega)})$.*

Proof: To find the heaviest $K_{2,k}$ we simply need to find, for any two vertices i, j, the k largest weighted vertices $v_1, \ldots, v_k$ so that each v_i is a common neighbor of i and j. As in Lemma 3.1, this reduces to finding the last k maximum witnesses of a 0-1 matrix product. A simple modification of the algorithm in Theorem 1.2 achieves this goal in the same running time (recall that k is fixed). $\square$

4   Heaviest H-subgraphs of real edge-weighted graphs

Given a vertex-colored graph G with n vertices, an H-subgraph of G is called *colorful* if each vertex of H has a distinct color. The *color coding* method presented in [AYZ95] is based upon two important facts. The first one is that, in many cases, finding a colorful H-subgraph is easier than finding an H-subgraph in an uncolored graph. The second one is that in a random vertex coloring with k colors, an H-subgraph with k vertices becomes colorful with probability $k!/k^k > e^{-k}$

and, furthermore, there is a derandomization technique that constructs a family of not too many colorings, so that each H-subgraph is colorful in at least one of the colorings. The derandomization technique, described in [AYZ95], constructs a family of colorings of size $O(\log n)$ whenever k is fixed. It is based upon a construction of *k-perfect hash functions* given in [FKS84] and in [SS90], and constructions of small probability spaces that admit almost ℓ-wise independent random variables [NN90]. The size of the constructed family of colorings is only $O(\log n)$ where k is fixed.

By the color coding method, in order to prove Theorem 1.4, it suffices to prove that, *given* a coloring of the vertices of the graph with k colors, a colorful cycle of length k of maximum weight (if one exists) can be found in $O(m^{2-1/\lceil k/2 \rceil})$ time.

Proof of Theorem 1.4: Assume that the vertices of G are colored with the colors $1, \ldots, k$. We first show that for each vertex u, a maximum weight colorful cycle of length k that passes through u can be found in $O(m)$ time. For a permutation π of $1, \ldots, k$, we show that a maximum weight cycle of the form $u = v_1, v_2, \ldots, v_k$ in which the color of v_i is $\pi(i)$ can be found in $O(m)$ time. Without loss of generality, assume π is the identity. For $j = 2, \ldots, k$ let V_j be the set of vertices whose color is j so that there is a path from u to $v \in V_j$ colored consecutively by the colors $1, \ldots, j$. Let $S(v)$ be the set of vertices of such a path with maximum possible weight. Denote this weight by $w(v)$. Clearly, V_j can be created from V_{j-1} in $O(m)$ time by examining the neighbors of each $v \in V_{j-1}$ colored with j. Now, let $w_u = \min_{v \in v_k} w(v) + w(v, u)$. Thus, w_u is the maximum weight of a cycle passing through u, of the desired form, and a cycle with this weight can be retrieved as well.

We prove the theorem when k is even. The odd case is similar. Let $\Delta = m^{2/k}$. There are at most $2m/\Delta = O(m^{1-2/k})$ vertices with degree at least Δ. For each vertex u with degree at least Δ we find a maximum weight colorful cycle of length k that passes through u. This can be done in $O(m^{2-2/k})$ time. It now suffices to find a maximum weight colorful cycle of length k in the subgraph G' of G induced by the vertices with maximum degree less than Δ. Consider a permutation π of $1, \ldots, k$. For a pair of vertices x, y, let S_1 be the set of all paths of length $k/2$ colored consecutively by $\pi(1), \ldots, \pi(k/2), \pi(k/2+1)$. There are at most $m\Delta^{k/2-1} = m^{2-2/k}$ such paths and they can be found using the greedy algorithm in $O(m^{2-2/k})$ time. Similarly, let S_2 be the set of all paths of length $k/2$ colored consecutively by $\pi(k/2+1), \ldots, \pi(k), \pi(1)$. If u, v are endpoints of at least one path in S_1 then let $f_1(\{u, v\})$ be the maximum weight of such a path. Similarly define $f_2(\{u, v\})$. We can therefore find, in $O(m^{2-2/k})$ a pair u, v (if one exists) so that $f_1(\{u, v\}) + f_2(\{u, v\})$ is maximized. By performing this procedure for each permutation, we find a maximum weight colorful cycle of length k in G'. $\qquad\square$

The definition of distance products, mentioned in the previous section, carries over to rectangular matrices. Let $A = A_{n_1 \times n_2}$ and $B = B_{n_2 \times n_3}$ be two matrices with entries in $\Re \cup \infty$. In this case, the *distance product* $C = A \star B$ is an $n_1 \times n_3$ matrix with $C[i, j] = \min_{k=1\ldots,n_2} A[i, k] + B[k, j]$. By partitioning the matrices into blocks it is obvious that Chan's algorithm for distance products of square matrices can be used to compute the distance product of an $n_1 \times n_2$ matrix and an $n_2 \times n_3$ matrix in $O(n_1 n_2 n_3 / \log \min\{n_1, n_2, n_3\})$ time. The analogous *MAX* version of distance products (namely, replacing MIN with MAX in the definition, and allowing $-\infty$) can be used to solve the

MAX H-SUBGRAPH problem in edge weighted graphs.

Proof of Theorem 1.5: We prove the theorem for $H = K_h$. The proof for other induced H-subgraphs is essentially the same. Partition h into a sum of three positive integers $a + b + c = h$. Let S_a be the set of all K_a-subgraphs of G. Notice that $|S_a| < n^a$ and that each $U \in S_a$ is an a-set. Similarly define S_b and S_c. We define A to be the matrix whose rows are indexed by S_a and whose columns are indexed by S_b. The entry $A[U, U']$ is defined to be $-\infty$ if $U \cup U'$ does not induce a K_{a+b}, otherwise it is defined to be the sum of the weights of the edges induced by $U \cup U'$. We define B to be the matrix whose rows are indexed by S_b and whose columns are indexed by S_c. The entry $A[U, U']$ is defined to by $-\infty$ if $U \cup U'$ does not induce a K_{b+c}, otherwise, it is defined to be the sum of the weights of the edges induced by $U \cup U'$ with *at least* one endpoint in U'. Notice the difference in the definitions of A and B. Let $C = A \star B$. The time to compute C using Chan's algorithm is $O(n^h / \log n)$. Now, for each $U \in S_a$ and $U' \in S_c$ so that $U \cup U'$ induces a K_{a+c}, let $w(U, U')$ be the sum of the weights of the edges with one endpoint in U and the other in U' plus the value of $C[U, U']$. If $w(U, U')$ is finite then it is the weight of the heaviest K_h that contains $U \cup U'$, otherwise no K_h contains $U \cup U'$. $\square$

The weighted DENSE k-SUBGRAPH problem (see, e.g., [FKP01]) is to find a k-vertex subgraph with maximum total edge weight. A simple modification of the algorithm of Theorem 1.5 solves this problem in $O(n^k / \log n)$ time. To our knowledge, this is the first non-trivial algorithm for this problem. Note that the maximum total weight of a k-subgraph can potentially be much larger than a k-clique's total weight.

Proof of Theorem 1.6: We use the color coding method, and an idea similar to Lemma 3.2 in [AYZ95]. Given a coloring of the vertices with k colors, it suffices to show how to find the heaviest colorful path of length $k - 1$ connecting any pair of vertices in $2^{O(k)} n^3 / \log n$ time. It will be convenient to assume that k is a power of two, and use recursion. Let C_1 be a set of $k/2$ distinct colors, and let C_2 be the complementary set of colors. Let V_i be the set of vertices colored by colors from C_i for $i = 1, 2$. Let G_i be the subgraph induced by V_i. Recursively find, for each pair of vertices in G_i, the maximum weight colorful path of length $k/2 - 1$. We record this information in matrices A_1, A_2, where the rows and columns of A_i are indexed by V_i. Let B be the matrix whose rows are indexed by V_1 and whose columns are indexed by V_2 where $B[u, v] = w(u, v)$. The max-distance product $D_{C_1, C_2} = (A_1 \star B) \star A_2$ gives, for each pair of vertices of G, all heaviest paths of length $k - 1$ where the first $k/2$ vertices are colored by colors from C_1 and the last $k/2$ vertices are colored by colors from C_2. By considering all $\binom{k}{k/2} < 2^k$ possible choices for (C_1, C_2), and computing D_{C_1, C_2} for each choice, we can obtain an $n \times n$ matrix D where $D[u, v]$ is the heaviest colorful path of length $k - 1$ between u and v. The number of distance products computed using this approach satisfies the recurrence $t(k) \leq 2^k t(k/2)$. Thus, the overall running time is $2^{O(k)} n^3 / \log n$. $\square$

The proof of Theorem 1.6 shows that, as long as $k = o(\log \log n)$, a cycle with k vertices and maximum weight can be found, with high probability, in $o(n^3)$ time. We note once again that all of our algorithms also apply to the *MIN* version of the problems. The previous best known algorithm for finding a minimum weight cycle of length k, in real weighted graphs, has running

time $O(k!n^3 2^k)$ [PV91].

5   Monochromatic and rainbow H-subgraphs

Proof of Theorem 1.7: Assume that H has t edges. The problem of finding a rainbow H-subgraph in G can be reduced, at a small cost, to the problem of finding a rainbow H-subgraph in another edge-coloring of G where the number of colors used is only t. Assume C is the set of colors used in G and consider a function $f : C \to \{1, \ldots, t\}$. This defines a new edge-coloring of G. Clearly, if an H-subgraph is not rainbow in the original coloring then it is also not rainbow in the new coloring. If f is constructed at random, a rainbow H-subgraph in the original coloring is also rainbow in the new coloring with probability $t!/t^t > e^{-t}$. As in the color coding method, this method can be derandomized by constructing $O(\log m) = O(\log n)$ colorings with only t colors used in each of them, so that if H is originally rainbow, it will also be rainbow in one of the constructed colorings.

We may now assume that $c : E \to \{1, \ldots, t\}$ and show how to find a rainbow H-subgraph if it exists, in $O(n^{\omega k + j})$ time. We shall assume that $H = K_h$ and $h = 3k + j$ where $j \in \{0, 1, 2\}$. The proof for other types of subgraphs is similar. By our assumption, $t = \binom{h}{2}$. Consider a partition of $\{1, \ldots, t\}$ into 6 parts $C_1, C_2, C_3, C_4, C_5, C_6$. The respective sizes are $|C_1| = |C_2| = \binom{k}{2}$, $|C_3| = \binom{k+j}{2}$, $|C_4| = k^2$, $|C_5| = |C_6| = k(k+j)$. Notice that there are $2^{O(t)}$ choices for the partition. For each partition we construct two Boolean matrices A and B that are defined as follows. The rows of A are indexed by all the rainbow K_k subgraphs of G that use the colors from C_1. The columns are indexed by all the rainbow K_k subgraphs of G that use the colors from C_2. We define $A[X, Y] = 1$ if $X \cap Y = \emptyset$ and the bipartite subgraph induced by the parts X and Y is complete, rainbow, and uses the colors from C_4, otherwise $A[X, Y] = 0$. The rows of B are indexed exactly in the same order as the columns of A. Namely, by all the rainbow K_k subgraphs of G that use the colors from C_2. The columns are indexed by all the rainbow K_{k+j} subgraphs of G that use the colors from C_3. We define $B[X, Y] = 1$ if $X \cap Y = \emptyset$ and the bipartite subgraph induced by the parts X and Y is complete, rainbow, and uses the colors from C_5, otherwise $A[X, Y] = 0$. The Boolean product $C = AB$ can be performed in $O(n^{\omega k + j})$ time using fast matrix multiplication. Now, if $C[X, Y] = 1$ and $X \cap Y = \emptyset$ and the bipartite subgraph of G induced by the parts X and Y is complete, rainbow, and uses colors from C_6 then we must have that $X \cup Y$ is contained in a rainbow K_h subgraph of G. By considering all possible partitions we are guaranteed not to miss a single rainbow K_h-subgraph of G. $\qquad\square$

Proof of Theorem 1.8: If H is a star, the theorem is trivial. Next, assume that H is not a star and has at least five vertices. Thus, H has two independent edges and at least one additional vertex. Put $h = 3k + j$ and consider a labeling of the vertices of H with $\{1, \ldots, h\}$ so that the following holds. If we partition $\{1, \ldots, h\}$ into three consecutive parts, as equally as possible, then the subgraph of H induced by the first part contains an edge e_1 and the subgraph induced by the second part contains an edge e_2. Thus, e.g., if H is the 5-cycle $(1, 2, 3, 4, 5)$ a plausible partition is $\{1, 2\}\{3, 4\}\{5\}$, $e_1 = (1, 2)$ and $e_2 = (3, 4)$. Denote by H_1, H_2, and H_3, the labeled subgraphs of H

induced by each of the parts and denote their respective sizes by h_1, h_1, h_3. Thus, if $j = 0$ we must have $h_1 = h_2 = h_3 = k$, if $j = 1$ we can assume $h_1 = k + 1$ and $h_2 = h_3 = k$ and if $j = 2$ we can assume $h_1 = h_2 = k + 1$ and $h_3 = k$.

We create a Boolean matrix A as follows. The rows of A are indexed by all the *ordered* h_1-tuples of vertices and the columns by all the *ordered* h_2-tuples. We put $A[X, Y] = 1$ if $X \cap Y = \emptyset$ and the mapping that assigns the i'th vertex of X to i and the ℓ'th vertex of Y to $h_1 + \ell$ corresponds to a monochromatic labeled copy of $H_1 \cup H_2$. In particular, note that the edge mapped to e_1 has the same color as the edge mapped to e_2. We create a Boolean matrix B as follows. The rows of B are indexed exactly like the columns of A. The columns of B are indexed by all the ordered h_3-tuples. We put $B[X, Y] = 1$ if $X \cap Y = \emptyset$ and the mapping that assigns the i'th vertex of X to $h_1 + i$ and the ℓ'th vertex of Y to $h_1 + h_2 + \ell$ corresponds to a monochromatic labeled copy of $H_2 \cup H_3$. Let $C = AB$ be the Boolean product. Suppose that $C[X, Y] = 1$ and suppose also that $X \cup Y$ corresponds to a monochromatic labeled copy of $H_1 \cup H_3$. Let Z satisfy $A[X, Z] = 1$ and $B[Z, Y] = 1$. Then we must have that $X \cup Y \cup Z$ corresponds to a monochromatic copy of H. This is because the color of each mapped edge is either that of the edge of G mapped to e_1 or that of the edge of G mapped to e_2 but these two are also colored the same. Notice also that if there is a monochromatic H-subgraph, it would be captured by our algorithm. Since the time needed to compute C is $O(n^{\omega k + j})$, the result follows.

Consider next the case $h = 4$. If $H \neq K_4$ then we can assume that H is labeled by $\{1, 2, 3, 4\}$ so that $(1, 4)$ is *not* an edge and $(2, 3)$ is an edge. Thus, the same algorithm described above using the partition $\{1\}\{2, 3\}\{4\}$ yields an $O(n^{\omega+1})$ time algorithm for detecting a monochromatic H. If $H = K_4$ the algorithm is slightly different. For each $v \in G$, let $S_1, \ldots, S_t$ be a partition of the neighbors of v so that $x, y \in S_i$ if and only if $c(v, x) = c(v, y)$. Searching for a triangle in the subgraph induced by S_i all of whose edges are colored by a given specific color has the same complexity as searching for an uncolored triangle in a graph, and hence can be done in $O(|S_i|^\omega)$ time. Thus, in $O(\sum_{i=1}^{t} |S_i|^\omega) \leq O(n^\omega)$ we can find a monochromatic triangle containing v, if it exists. Performing this procedure for each $v \in V$ gives the desired $O(n^{\omega+1})$ time algorithm.

The only remaining case is $H = K_3$. Let E_i be the set of edges of G colored with i. We say that i is *heavily used* if $|E_i| \geq n^{(\omega+1)/2}$. Clearly, the number of colors heavily used is at most $O(n^{2-(\omega+1)/2})$. For each heavily used color i we can decide, in $O(n^\omega)$ time, whether there is a monochromatic triangle colored with i. The overall running time is, therefore $O(n^{\omega+2-(\omega+1)/2}) = O(n^{(3+\omega)/2})$. For each color i that is not heavily used, we can decide in $O(|E_i|^{2\omega/(\omega+1)})$ time whether there is a monochromatic triangle colored with i using the algorithm from [AYZ97]. The overall running time is maximized if $|E_i| = \Theta(n^{(\omega+1)/2})$ and when there are $\Theta(n^{2-(\omega+1)/2})$ such colors. In this extremal case the running time is still only $O(n^{((\omega+1)/2)(2\omega/(\omega+1))+2-(\omega+1)/2}) = O(n^{(3+\omega)/2})$. $\square$

6 Concluding remarks and open problems

We presented several algorithms for MAX H-SUBGRAPH in both real vertex weighted or real edge weighted graphs, and results for the related problem of finding monochromatic or rainbow

H-subgraphs in edge-colored graphs. It may be possible to improve upon the running times of some of our algorithms. More specifically, we raise the following open problems.

(i) Can the exponent $t(\omega, 3)$ in Theorem 1.2 be improved? If so, this would immediately imply an improved algorithm for maximum witnesses.

(ii) Can the logarithmic factor in Theorem 1.4 be eliminated? We know from [AYZ97] that this is the case in the unweighted version of the problem. Can the logarithmic factor in Theorem 1.7 be eliminated?

(iii) Can monochromatic triangles be detected faster than the $O(n^{(3+\omega)/2})$ algorithm of Theorem 1.8? In particular, can they be detected in $O(n^\omega)$ time?

(iv) Is the dominance matrix for a set of n points in n dimensions computable in $\tilde{O}(n^\omega)$ time?

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